

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSPTADEG1625

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *
SESSION RESUMED IN FILE 'CAPLUS' AT 15:16:46 ON 07 JUN 2007
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| COST IN U.S. DOLLARS | SINCE FILE ENTRY | TOTAL SESSION |
|----------------------|---------------------|------------------|
| FULL ESTIMATED COST | 6.37 | 35.61 |

=> file caplus

| COST IN U.S. DOLLARS | SINCE FILE ENTRY | TOTAL SESSION |
|----------------------|---------------------|------------------|
| FULL ESTIMATED COST | 6.37 | 35.61 |

FILE 'CAPLUS' ENTERED AT 15:16:59 ON 07 JUN 2007
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FILE COVERS 1907 - 7 Jun 2007 VOL 146 ISS 24
FILE LAST UPDATED: 6 Jun 2007 (20070606/ED)

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=> s aird/au

L7 0 AIRD/AU

=> log off

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF
LOGOFF? (Y)/N/HOLD:n

=> s ruthenium

96839 RUTHENIUM
23 RUTHENIUMS
L8 96839 RUTHENIUM
(RUTHENIUM OR RUTHENIUMS)

=> s 18 and carmona/au

0 CARMONA/AU
L9 0 L8 AND CARMONA/AU

=> log hold

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

7.91

43.52

SESSION WILL BE HELD FOR 120 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 15:19:12 ON 07 JUN 2007

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSPTADEG1625

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * *

SESSION RESUMED IN FILE 'CAPLUS' AT 15:22:27 ON 07 JUN 2007

FILE 'CAPLUS' ENTERED AT 15:22:27 ON 07 JUN 2007

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COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

7.91

43.52

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

7.91

43.52

FILE 'CAPLUS' ENTERED AT 15:22:44 ON 07 JUN 2007

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FILE COVERS 1907 - 7 Jun 2007 VOL 146 ISS 24

FILE LAST UPDATED: 6 Jun 2007 (20070606/ED)

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=> e carmona d/au

E1 5 CARMONA CYNTHIA/AU

E2 3 CARMONA CYNTHIA L/AU

E3 25 --> CARMONA D/AU

E4 81 CARMONA DANIEL/AU

E5 1 CARMONA DAVID/AU

E6 1 CARMONA DE CATRIL A/AU

E7 1 CARMONA DE GARCIA C A/AU

| | | |
|-----|----|---------------------------|
| E8 | 1 | CARMONA DIAZ ELIZABETH/AU |
| E9 | 1 | CARMONA DORIS/AU |
| E10 | 55 | CARMONA E/AU |
| E11 | 6 | CARMONA E C/AU |
| E12 | 1 | CARMONA E FERREIRA R/AU |

o

=> s carmona d?/au and ruthenium?

111 CARMONA D?/AU

96878 RUTHENIUM?

L10 . 34 CARMONA D?/AU AND RUTHENIUM?

=> l10 and pyrazolyl

L10 IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system.
For a list of commands available to you in the current file, enter
"HELP COMMANDS" at an arrow prompt (=>).

=> s l10 and pyrazolyl

6857 PYRAZOLYL

8 PYRAZOLYLS

6859 PYRAZOLYL

(PYRAZOLYL OR PYRAZOLYLS)

L11 3 L10 AND PYRAZOLYL

=> d l11 1-3 abs ibib hitstr

L11 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2007 ACS on STN

AB The reaction of the metallo-ligand [Ru(η 6-p-cymene) (pz)₂ (Hpz)] with the Pt complex [{Pt(Ime)₃]₄] affords mixts. of heterodinuclear [(η 6-p-cymene)Ru(μ -pz)₃PtMe₃] (1) and [(η 6-p-cymene)Ru(μ -pz)₂(μ -I)PtMe₃] (2). The reaction of the Ir derivative [Ir(η 5-C₅Me₅) (pz)₂ (Hpz)] with [{Pt(Ime)₃]₄] gives [(η 5-C₅Me₅)Ir(μ -pz)₂(μ -I)PtMe₃] (3). Both [Ru(η 6-p-cymene) (pz)₂ (Hpz)] and [Ir(η 5-C₅Me₅) (pz)₂ (Hpz)] react with [{Pt(Ime)₃]₄] in the presence of NaOH yielding 1 and [(η 5-C₅Me₅)Ir(μ -pz)₃PtMe₃] (4), resp. While [Ru(η 6-p-cymene) (pz)₂ (Hpz)] reacts with [PtBr₂Me₂Sx] to give mixts. of [(η 6-p-cymene)Ru-(μ -pz)₃PtBrMe₂] (5) and [(η 6-p-cymene)Ru(μ -pz)₂(μ -Br)PtBrMe₂] (6), the reaction of [Ir(η 5-C₅Me₅) (pz)₂ (Hpz)] with [PtBr₂Me₂Sx] gives [(η 5-C₅Me₅)Ir(μ -pz)₂(μ -Br)PtBrMe₂] (7) as the sole product. All species were characterized in solution by 1H-NMR spectroscopy. The crystal structure of complex 4 was determined by single-crystal x-ray diffraction.

ACCESSION NUMBER: 2000:695725 CAPLUS

DOCUMENT NUMBER: 134:5034

TITLE: Synthesis and characterization of heterodinuclear RuPt and IrPt complexes containing pyrazolate bridging ligands. Crystal structure of [(η 5-C₅Me₅)Ir(μ -pz)₃PtMe₃] (pz = pyrazolate)

AUTHOR(S): Contreras, Raul; Valderrama, Mauricio; Orellana, Esteban M.; Boys, Daphne; Carmona, Daniel;

CORPORATE SOURCE: Oro, Luis A.; Lamata, M. Pilar; Ferrer, Joaquina
Departamento de Quimica Inorganica, Facultad de Quimica, Pontificia Universidad Catolica de Chile, Santiago, 22, Chile

SOURCE: Journal of Organometallic Chemistry (2000), 606(2), 197-202

CODEN: JORCAI; ISSN: 0022-328X

PUBLISHER: Elsevier Science S.A.

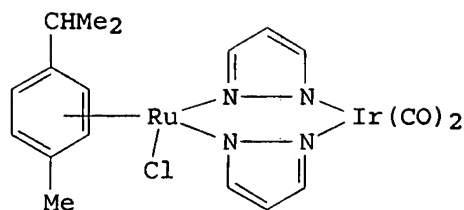
DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 134:5034

REFERENCE COUNT: 36 THERE ARE 36 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L11 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2007 ACS on STN



I

AB Dinuclear iridium-ruthenium complex I (M = Ir), prepared from (η^6 -p-cymol)RuCl(μ -pz)Ir(COD) (pz = pyrazolyl, COD = 1,5-cyclooctadiene) and carbon monoxide, reacted with NaBr or NaI to give the corresponding halo complex [(η^6 -p-cymol)Ru(μ -pz)2IrX(CO)2] (X = Br, iodo). Complexes I (M = Ir, Rh) were characterized by x-ray crystallog.

ACCESSION NUMBER: 1991:656346 CAPLUS

DOCUMENT NUMBER: 115:256346

TITLE: Reversible isomerization of the dinuclear complex [(η^6 -p-Cymol)RuCl(μ -Pyrazolyl)2Ir(CO)2] with formation of a ruthenium-iridium bond

AUTHOR(S): Carmona, Daniel; Ferrer, Joaquina; Mendoza, Ana; Lahoz, Fernando J.; Reyes, Josefa; Oro, Luis A.

CORPORATE SOURCE: Inst. Cienc. Mater. Aragon, Univ. Zaragoza, Zaragoza, E-50009, Spain

SOURCE: Angewandte Chemie (1991), 103(9), 1192-4 (See also Angew. Chem., Int. Ed. Engl., 1991, 30(9), 1171-3) CODEN: ANCEAD; ISSN: 0044-8249

DOCUMENT TYPE: Journal

LANGUAGE: English

L11 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2007 ACS on STN

AB The synthesis and NMR spectra (^1H and ^{13}C) are reported for 24 p-cymeneruthenium complexes belonging to one of the following families: [Ru(MeC6H4CHMe2-p)(acac)X], [Ru(MeC6H4CHMe2-p)(acac)L]BF4, [Ru(MeC6H4CHMe2-p)ClL2]BF4, and [Ru(MeC6H4CHMe2-p)L3][BF4]2, [Ru(MeC6H4CHMe2-p)XL2]BF4, and [Ru(MeC6H4CHMe2-p)X2L] where X = Br, I, N3, pz, mpz, dmpz, or idz, and L = pyridine, PPh3, CNCMe3, P(OMe)3, Hpz (pyrazole), Hmpz (3-methylpyrazole), Hdmpz (3,5-dimethylpyrazole), and Hidz (indazole) for some complexes and only azoles (pyrazoles and indazole) for the remaining ones. Crystals of [Ru(MeC6H4CHMe2-p)(pz)(Hpz)2]BF4 are obtained and the structure was determined by x-ray diffraction. There are 2 crystallog. units, each having an intramol. hydrogen bond between a pyrazole and a pyrazolate ring, and another between the other pyrazole ligand and the BF4 anion. The NMR data (δ and J) of the azole complexes were carefully determined and are thoroughly discussed.

ACCESSION NUMBER: 1990:478665 CAPLUS

DOCUMENT NUMBER: 113:78665

TITLE: Synthesis, x-ray structure, and nuclear magnetic resonance proton and carbon-13 studies of ruthenium(II) complexes containing pyrazolyl ligands

AUTHOR(S): Carmona, Daniel; Ferrer, Joaquina; Oro, Luis A.; Apreda, Maria C.; Foces-Foces, Concepcion; Cano, Felix H.; Elguero, Jose; Luisa Jimeno, Maria

CORPORATE SOURCE: Inst. Cienc. Mater. Aragon, Univ. Zaragoza, Zaragoza, 50009, Spain

SOURCE: Journal of the Chemical Society, Dalton Transactions: Inorganic Chemistry (1972-1999) (1990), (4), 1463-76 CODEN: JCDTBI; ISSN: 0300-9246

DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 113:78665

=> analyze l11 3

ENTER DISPLAY CODE (TI) OR ?:end

=> analyze

ENTER ANSWER SET OR ANALYZE L# OR (L11):l11

ENTER ANSWER NUMBER OR RANGE (1-):3

ENTER DISPLAY CODE (TI) OR ?:ab

L12 ANALYZE L11 3 AB : 85 TERMS

=> d doc

L12 ANALYZE L11 3 AB : 85 TERMS

| TERM # | # OCC | # DOC | % DOC | AB |
|--------|-------|-------|-------|----|
|--------|-------|-------|-------|----|

| | | | | |
|----|---|---|--------|-------------|
| 1 | 9 | 1 | 100.00 | P |
| 2 | 7 | 1 | 100.00 | MEC6H4CHME2 |
| 3 | 7 | 1 | 100.00 | RU |
| 4 | 6 | 1 | 100.00 | BF4 |
| 5 | 4 | 1 | 100.00 | ARE |
| 6 | 3 | 1 | 100.00 | COMPLEXES |
| 7 | 3 | 1 | 100.00 | PYRAZOLE |
| 8 | 3 | 1 | 100.00 | X |
| 9 | 3 | 1 | 100.00 | 2 |
| 10 | 3 | 1 | 100.00 | 3 |

75 MORE TERMS WITH A DOCUMENT COUNT OF 1

=> analyze

ENTER ANSWER SET OR ANALYZE L# OR (L12):l11

ENTER ANSWER NUMBER OR RANGE (1-):3

ENTER DISPLAY CODE (TI) OR ?:cc

L13 ANALYZE L11 3 CC : 1 TERM

=> d doc

L13 ANALYZE L11 3 CC : 1 TERM

| TERM # | # OCC | # DOC | % DOC | CC |
|--------|-------|-------|-------|----|
|--------|-------|-------|-------|----|

| | | | | |
|---|---|---|--------|-------|
| 1 | 1 | 1 | 100.00 | 29-13 |
|---|---|---|--------|-------|

***** END OF L13***

=> analyze

ENTER ANSWER SET OR ANALYZE L# OR (L13):l11

ENTER ANSWER NUMBER OR RANGE (1-):3

ENTER DISPLAY CODE (TI) OR ?:rn

L14 ANALYZE L11 3 RN : 34 TERMS

=> d doc

L14 ANALYZE L11 3 RN : 34 TERMS

| TERM # | # OCC | # DOC | % DOC | RN |
|--------|-------|-------|-------|----|
|--------|-------|-------|-------|----|

| | | | | |
|---|---|---|--------|-------------|
| 1 | 1 | 1 | 100.00 | 121-45-9 |
| 2 | 1 | 1 | 100.00 | 128628-32-0 |
| 3 | 1 | 1 | 100.00 | 128628-33-1 |
| 4 | 1 | 1 | 100.00 | 128628-34-2 |
| 5 | 1 | 1 | 100.00 | 128628-35-3 |
| 6 | 1 | 1 | 100.00 | 128628-36-4 |
| 7 | 1 | 1 | 100.00 | 128628-37-5 |

8 1 1 100.00 128628-38-6
9 1 1 100.00 128628-40-0
10 1 1 100.00 128628-42-2
24 MORE TERMS WITH A DOCUMENT COUNT OF 1

=> FILE REG

| COST IN U.S. DOLLARS | SINCE FILE ENTRY | TOTAL SESSION |
|--|------------------|---------------|
| FULL ESTIMATED COST | 64.10 | 107.62 |
| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE ENTRY | TOTAL SESSION |
| CA SUBSCRIBER PRICE | -2.34 | -2.34 |

FILE 'REGISTRY' ENTERED AT 15:43:02 ON 07 JUN 2007
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DICTIONARY FILE UPDATES: 6 JUN 2007 HIGHEST RN 936692-95-4

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experimental property data in the original document. For information
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<http://www.cas.org/support/stngen/stndoc/properties.html>

=> STR 121-45-9

:END

L15 STRUCTURE CREATED

=> S L15 EXA SAM

SAMPLE SEARCH INITIATED 15:43:06 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L16 0 SEA EXA SAM L15

=>

=>

NO ANSWERS WERE FOUND

=> FILE REG

| | | |
|--|------------|---------|
| COST IN U.S. DOLLARS | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| FULL ESTIMATED COST | 0.90 | 108.52 |
| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| CA SUBSCRIBER PRICE | 0.00 | -2.34 |

FILE 'REGISTRY' ENTERED AT 15:43:57 ON 07 JUN 2007
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on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=> STR 121-45-9

:END

L17 STRUCTURE CREATED

=> S L17 EXA FUL

FULL SEARCH INITIATED 15:44:00 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 59 TO ITERATE

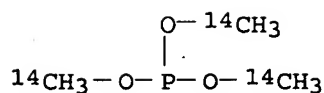
100.0% PROCESSED 59 ITERATIONS 9 ANSWERS
SEARCH TIME: 00.00.01

L18 9 SEA EXA FUL L17

=>

=> D SCAN

L18 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN
IN Methanol-14C, phosphite (3:1) (9CI)
MF C3 H9 O3 P



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> d l14 hitstr

L14 ANALYZE L11 3 RN : 34 TERMS

NO TERMS MATCHED THE DISPLAY CRITERIA

=> FILE REG

| | | |
|--|------------|---------|
| COST IN U.S. DOLLARS | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| FULL ESTIMATED COST | 60.95 | 169.47 |
| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE | TOTAL |
| | ENTRY | SESSION |
| CA SUBSCRIBER PRICE | 0.00 | -2.34 |

FILE 'REGISTRY' ENTERED AT 15:48:22 ON 07 JUN 2007
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 DICTIONARY FILE UPDATES: 6 JUN 2007 HIGHEST RN 936692-95-4

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 conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
 predicted properties as well as tags indicating availability of
 experimental property data in the original document. For information
 on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=> STR 128628-32-0

:END

L19 STRUCTURE CREATED

=> S L19 EXA FUL

FULL SEARCH INITIATED 15:48:26 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 1 TO ITERATE

100.0% PROCESSED 1 ITERATIONS 1 ANSWERS
 SEARCH TIME: 00.00.01

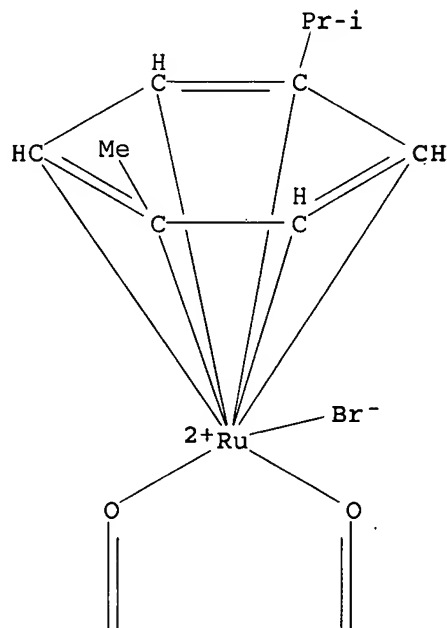
L20 1 SEA EXA FUL L19

=>

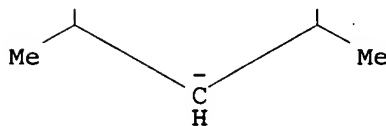
=> D SCAN

L20 1 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN
IN Ruthenium, bromo[(1,2,3,4,5,6-η)-1-methyl-4-(1-methylethyl)benzene] (2,4-pentanedionato-O,O')- (9CI)
MF C15 H21 Br O2 Ru
CI CCS

PAGE 1-A



PAGE 2-A



ALL ANSWERS HAVE BEEN SCANNED

=> s l19 and (anticancer or antitumor or antitumour)
COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID
The query entered contains both search terms created by structure-building or screen commands and text search terms. L#s created via the STRUCTURE or SCREEN commands must be searched in the structures files separately from text terms or profiles. The L# answer sets from structure searches can be used in crossover searches and can be combined with text terms.

=> s l20 and (anticancer or antitumor or antitumour)

74 ANTICANCER
209 ANTITUMOR
0 ANTITUMOUR
L21 0 L20 AND (ANTICANCER OR ANTITUMOR OR ANTITUMOUR)

=> log hold

COST IN U.S. DOLLARS

| SINCE FILE | TOTAL |
|------------|---------|
| ENTRY | SESSION |
| 78.05 | 247.52 |

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

| SINCE FILE | TOTAL |
|------------|---------|
| ENTRY | SESSION |
| 0.00 | -2.34 |

CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 120 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 15:55:21 ON 07 JUN 2007